

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN013024\
 Data File : BN029414.D
 Acq On : 30 Jan 2024 14:46
 Operator : MA/JU
 Sample : 03572-02
 Misc : LOQ-MDL-SOIL-0.2NG
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 LOQ-SOIL-02-QT3-2023

Quant Time: Jan 30 18:17:24 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN012924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 30 03:38:24 2024
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 01/31/2024
 Supervised By :Sohil Jodhani 02/02/2024

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.898	152	3783	0.400	ng	0.00
7) Naphthalene-d8	10.701	136	10046	0.400	ng	0.00
13) Acenaphthene-d10	14.532	164	4456	0.400	ng	-0.01
19) Phenanthrene-d10	17.293	188	8830	0.400	ng	0.00
29) Chrysene-d12	21.492	240	6043	0.400	ng	# 0.00
35) Perylene-d12	23.879	264	6291	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.464	112	3265	0.359	ng	0.00
5) Phenol-d6	7.060	99	3832	0.365	ng	0.00
8) Nitrobenzene-d5	9.067	82	2883	0.345	ng	0.00
11) 2-Methylnaphthalene-d10	12.291	152	7090	0.518	ng	0.00
14) 2,4,6-Tribromophenol	16.026	330	470	0.329	ng	-0.01
15) 2-Fluorobiphenyl	13.164	172	6587	0.344	ng	-0.01
27) Fluoranthene-d10	19.327	212	12080	0.565	ng	0.00
31) Terphenyl-d14	19.935	244	4997	0.333	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.377	88	1133	0.213	ng	98
3) n-Nitrosodimethylamine	3.716	42	728	0.168	ng	99
6) bis(2-Chloroethyl)ether	7.334	93	1699	0.170	ng	96
9) Naphthalene	10.744	128	5114	0.179	ng	98
10) Hexachlorobutadiene	11.021	225	961	0.177	ng	# 98
12) 2-Methylnaphthalene	12.363	142	2825	0.168	ng	97
16) Acenaphthylene	14.254	152	3959	0.187	ng	99
17) Acenaphthene	14.596	154	2505	0.176	ng	98
18) Fluorene	15.580	166	3144	0.176	ng	97
20) 4,6-Dinitro-2-methylph...	15.667	198	270	0.208	ng	# 72
21) 4-Bromophenyl-phenylether	16.486	248	909	0.173	ng	91
22) Hexachlorobenzene	16.585	284	1148	0.178	ng	99
23) Atrazine	16.759	200	677	0.181	ng	# 89
24) Pentachlorophenol	16.933	266	357	0.175	ng	98
25) Phenanthrene	17.330	178	4665	0.179	ng	99
26) Anthracene	17.429	178	3709	0.167	ng	97
28) Fluoranthene	19.354	202	4913	0.167	ng	100
30) Pyrene	19.721	202	4937	0.177	ng	99
32) Benzo(a)anthracene	21.474	228	3448	0.167	ng	99
33) Chrysene	21.528	228	4172	0.178	ng	99
34) Bis(2-ethylhexyl)phtha...	21.429	149	2419	0.172	ng	99
36) Indeno(1,2,3-cd)pyrene	26.341	276	4140	0.164	ng	# 81
37) Benzo(b)fluoranthene	23.154	252	3622	0.161	ng	# 89
38) Benzo(k)fluoranthene	23.204	252	4055m	0.172	ng	
39) Benzo(a)pyrene	23.771	252	3183	0.162	ng	# 83
40) Dibenzo(a,h)anthracene	26.376	278	3221	0.160	ng	# 92
41) Benzo(g,h,i)perylene	27.092	276	3989	0.163	ng	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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