

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041823\
 Data File : BN024972.D
 Acq On : 18 Apr 2023 11:08
 Operator : CG/JU
 Sample : 02357-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 P2-S2-230413

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 04/19/2023
 Supervised By :Jagrut Upadhyay 04/19/2023

Quant Time: Apr 18 11:43:16 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN040523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 18 05:29:26 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.933	152	11522	0.400	ng	0.00
7) Naphthalene-d8	10.739	136	34979	0.400	ng	#-0.01
13) Acenaphthene-d10	14.576	164	20110	0.400	ng	0.00
19) Phenanthrene-d10	17.323	188	47960	0.400	ng	0.00
29) Chrysene-d12	21.522	240	47335	0.400	ng	# 0.00
35) Perylene-d12	23.975	264	50772	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.484	112	8837	0.332	ng	0.00
5) Phenol-d6	7.095	99	11539	0.344	ng	0.00
8) Nitrobenzene-d5	9.095	82	8258	0.293	ng	-0.01
11) 2-Methylnaphthalene-d10	12.332	152	14058	0.303	ng	0.00
14) 2,4,6-Tribromophenol	16.070	330	3304	0.365	ng	0.00
15) 2-Fluorobiphenyl	13.197	172	21150	0.275	ng	-0.01
27) Fluoranthene-d10	19.355	212	33229	0.306	ng	0.00
31) Terphenyl-d14	19.949	244	29102	0.271	ng	0.00
Target Compounds						Qvalue
6) bis(2-Chloroethyl)ether	7.355	93	2346	0.069	ng	# 89
9) Naphthalene	10.792	128	100601	1.117	ng	99
12) 2-Methylnaphthalene	12.404	142	34149	0.545	ng	99
16) Acenaphthylene	14.298	152	140531	1.662	ng	99
17) Acenaphthene	14.640	154	29282	0.512	ng	99
18) Fluorene	15.624	166	108145	1.378	ng	98
25) Phenanthrene	17.360	178	1336526	9.598	ng	100
26) Anthracene	17.447	178	451000	3.715	ng	100
28) Fluoranthene	19.384	202	2262776	14.314	ng	99
30) Pyrene	19.751	202	1877749	11.160	ng	96
32) Benzo(a)anthracene	21.504	228	1079726	6.746	ng	97
33) Chrysene	21.558	228	962195	5.739	ng	98
34) Bis(2-ethylhexyl)phtha...	21.423	149	148232	1.724	ng	98
36) Indeno(1,2,3-cd)pyrene	26.528	276	497255	2.234	ng	95
37) Benzo(b)fluoranthene	23.227	252	1251420	6.499	ng	# 94
38) Benzo(k)fluoranthene	23.268	252	451901m	2.368	ng	
39) Benzo(a)pyrene	23.867	252	911795m	4.951	ng	
40) Dibenzo(a,h)anthracene	26.537	278	119123	0.680	ng	94
41) Benzo(g,h,i)perylene	27.311	276	490797	2.586	ng	99

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

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