

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041621\
 Data File : BN014319.D
 Acq On : 16 Apr 2021 20:47
 Operator : CG/JU
 Sample : M2061-02
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GWBR3-225-250-041521

Quant Time: Apr 17 02:53:42 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN041621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 16 16:29:41 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.700	152	7593	0.40	ng	0.00
7) Naphthalene-d8	10.467	136	28629	0.40	ng	#-0.01
13) Acenaphthene-d10	14.334	164	15642	0.40	ng	0.00
19) Phenanthrene-d10	17.068	188	33565	0.40	ng	0.00
29) Chrysene-d12	21.271	240	34716	0.40	ng	0.00
36) Perylene-d12	23.496	264	29661	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.300	112	1983	0.12	ng	0.00
5) Phenol-d6	6.871	99	1433	0.07	ng	0.00
8) Nitrobenzene-d5	8.845	82	5789	0.32	ng	0.00
11) 2-Methylnaphthalene-d10	12.066	152	13120	0.30	ng	0.00
14) 2,4,6-Tribromophenol	15.818	330	2207	0.39	ng	0.00
15) 2-Fluorobiphenyl	12.951	172	20304	0.33	ng	0.00
27) Fluoranthene-d10	19.109	212	37006	0.40	ng	0.00
31) Terphenyl-d14	19.715	244	42849	0.59	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.255	88	104050	10.79	ng	93
9) Naphthalene	10.518	128	2761	0.04	ng	# 92
12) 2-Methylnaphthalene	12.137	142	2946	0.06	ng	97
24) Pentachlorophenol	16.727	266	134	0.02	ng	92
25) Phenanthrene	17.116	178	7944	0.08	ng	99
28) Fluoranthene	19.139	202	9038	0.09	ng	# 87
30) Pyrene	19.501	202	7445	0.08	ng	# 95
32) Benzo(a)anthracene	21.250	228	5657	0.06	ng	95
33) Chrysene	21.303	228	7917	0.07	ng	96
34) Bis(2-ethylhexyl)phtha...	21.186	149	13859	0.53	ng	100
35) Indeno(1,2,3-cd)pyrene	25.735	276	8474	0.06	ng	97
37) Benzo(b)fluoranthene	22.825	252	11756	0.09	ng	# 80
38) Benzo(k)fluoranthene	22.870	252	8711	0.07	ng	# 73
39) Benzo(a)pyrene	23.397	252	4724	0.05	ng	# 46
40) Dibenzo(a,h)anthracene	25.752	278	5771	0.05	ng	# 82
41) Benzo(g,h,i)perylene	26.419	276	7919	0.07	ng	# 91

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

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