

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031921\
 Data File : BN013986.D
 Acq On : 18 Mar 2021 17:31
 Operator : CG/JU
 Sample : M1714-01DL 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GWBR3-224-234-031721DL

Quant Time: Mar 18 18:00:56 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN022421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 18 15:58:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.708	152	6350	0.40	ng	0.00
7) Naphthalene-d8	10.492	136	24033	0.40	ng	# 0.00
13) Acenaphthene-d10	14.346	164	13476	0.40	ng	0.00
19) Phenanthrene-d10	17.092	188	30517	0.40	ng	# 0.00
29) Chrysene-d12	21.271	240	30903	0.40	ng	# 0.00
36) Perylene-d12	23.503	264	27457	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.300	112	459	0.03	ng	0.00
5) Phenol-d6	6.871	99	395	0.02	ng	0.00
8) Nitrobenzene-d5	8.857	82	847	0.05	ng	0.00
11) 2-Methylnaphthalene-d10	12.080	152	1950	0.05	ng	0.00
14) 2,4,6-Tribromophenol	15.831	330	309	0.06	ng	0.00
15) 2-Fluorobiphenyl	12.963	172	2897	0.06	ng	0.00
27) Fluoranthene-d10	19.119	212	5713	0.06	ng	0.00
31) Terphenyl-d14	19.720	244	6105	0.09	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.247	88	15330	2.00	ng	97
12) 2-Methylnaphthalene	12.155	142	907	0.02	ng	# 94
25) Phenanthrene	17.128	178	2403	0.03	ng	98
28) Fluoranthene	19.149	202	5961	0.06	ng	98
30) Pyrene	19.511	202	5336	0.06	ng	97
32) Benzo(a)anthracene	21.261	228	3547	0.04	ng	94
33) Chrysene	21.303	228	5139	0.06	ng	97
34) Bis(2-ethylhexyl)phtha...	21.197	149	83785	1.85	ng	99
35) Indeno(1,2,3-cd)pyrene	25.735	276	3554	0.03	ng	96
37) Benzo(b)fluoranthene	22.832	252	6207	0.07	ng	# 74
38) Benzo(k)fluoranthene	22.877	252	3171	0.04	ng	# 65
39) Benzo(a)pyrene	23.407	252	3588	0.05	ng	# 50
41) Benzo(g,h,i)perylene	26.419	276	3627	0.04	ng	92

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

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