

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE041221\
 Data File : BE101287.D
 Acq On : 12 Apr 2021 19:57
 Operator : CG/JU
 Sample : M1966-03DL 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MW-110RRDL

Quant Time: Apr 13 01:45:58 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-SIM-BE041221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 12 18:02:55 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.830	152	2510	0.40	ng	0.00
7) Naphthalene-d8	10.614	136	12596	0.40	ng	0.00
13) Acenaphthene-d10	14.442	164	10133	0.40	ng	0.00
19) Phenanthrene-d10	17.196	188	27886	0.40	ng	# 0.00
29) Chrysene-d12	21.353	240	38915	0.40	ng	0.00
36) Perylene-d12	23.841	264	27897	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.412	112	973	0.13	ng	-0.02
5) Phenol-d6	7.001	99	381	0.03	ng	0.00
8) Nitrobenzene-d5	8.978	82	1705	0.15	ng	0.00
11) 2-Methylnaphthalene-d10	12.195	152	3321	0.11	ng	0.00
14) 2,4,6-Tribromophenol	15.957	330	400	0.13	ng	0.00
15) 2-Fluorobiphenyl	13.073	172	4993	0.14	ng	0.00
27) Fluoranthene-d10	19.216	212	70437	0.14	ng	0.00
31) Terphenyl-d14	19.813	244	16813	0.18	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.362	88	431	0.10	ng	# 1
9) Naphthalene	10.666	128	476488	3.50	ng	100
12) 2-Methylnaphthalene	12.267	142	1644	0.07	ng	# 94
16) Acenaphthylene	14.168	152	3549	0.09	ng	# 87
17) Acenaphthene	14.500	154	5290	0.18	ng	98
18) Fluorene	15.503	166	8093	0.21	ng	# 97
24) Pentachlorophenol	16.864	266	146	0.34	ng	# 87
25) Phenanthrene	17.242	178	30060	0.37	ng	# 94
26) Anthracene	17.332	178	19172	0.26	ng	# 82
28) Fluoranthene	19.245	202	31492	0.30	ng	99
30) Pyrene	19.606	202	27533	0.22	ng	98
32) Benzo(a)anthracene	21.341	228	13825	0.11	ng	95
33) Chrysene	21.390	228	11585	0.09	ng	# 92
34) Bis(2-ethylhexyl)phtha...	21.269	149	14660	0.09	ng	100
37) Benzo(b)fluoranthene	23.078	252	9725	0.10	ng	# 88
38) Benzo(k)fluoranthene	23.128	252	7742	0.08	ng	# 85
39) Benzo(a)pyrene	23.727	252	6419	0.08	ng	# 83

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

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