

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040621\
 Data File : BP005210.D
 Acq On : 06 Apr 2021 22:49
 Operator : CG/JU
 Sample : M1909-02DL 10X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FS-SB-10-26-27DL

Quant Time: Apr 07 00:32:42 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 06 12:33:08 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.693	152	27010	20.00	ng	# 0.00
21) Naphthalene-d8	10.493	136	106254	20.00	ng	# 0.00
39) Acenaphthene-d10	14.351	164	64192	20.00	ng	0.00
64) Phenanthrene-d10	17.116	188	136044	20.00	ng	# 0.00
76) Chrysene-d12	21.210	240	136614	20.00	ng	0.00
86) Perylene-d12	23.492	264	140985	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.293	112	18794	10.71	ng	0.00
7) Phenol-d6	6.875	99	27726	11.03	ng	0.00
23) Nitrobenzene-d5	8.852	82	17837	7.20	ng	0.00
42) 2,4,6-Tribromophenol	15.851	330	6078	7.27	ng	0.00
45) 2-Fluorobiphenyl	12.963	172	38331	8.08	ng	0.00
79) Terphenyl-d14	19.686	244	57147	7.61	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.563	128	4902803	816.81	ng	99
37) 2-Methylnaphthalene	12.163	142	584919	134.96	ng	# 96
38) 1-Methylnaphthalene	12.381	142	298871	73.57	ng	97
46) 1,1'-Biphenyl	13.181	154	38486	7.86	ng	95
49) Acenaphthylene	14.075	152	19229	3.10	ng	# 90
71) Phenanthrene	17.157	178	47413	6.01	ng	99
75) Fluoranthene	19.134	202	47036	5.07	ng	97
78) Pyrene	19.486	202	45462	4.84	ng	97
81) Benzo(a)anthracene	21.198	228	21877	2.36	ng	96
83) Chrysene	21.245	228	20149	2.21	ng	# 93
88) Benzo(b)fluoranthene	22.804	252	21003	2.10	ng	# 92

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

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