

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112921\
 Data File : BP008137.D
 Acq On : 29 Nov 2021 12:35
 Operator : CG/JU
 Sample : M4815-02DL 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-5DL

Quant Time: Nov 29 12:59:49 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 25 01:13:46 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	105239	20.000	ng	-0.01
21) Naphthalene-d8	10.604	136	381429	20.000	ng	# 0.00
39) Acenaphthene-d10	14.451	164	237111	20.000	ng	0.00
64) Phenanthrene-d10	17.210	188	498750	20.000	ng	0.00
76) Chrysene-d12	21.310	240	587898	20.000	ng	0.00
86) Perylene-d12	23.704	264	657757	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	39566	6.053	ng	-0.01
7) Phenol-d6	6.987	99	31594	3.581	ng	-0.01
23) Nitrobenzene-d5	8.963	82	83823	9.993	ng	-0.01
42) 2,4,6-Tribromophenol	15.951	330	38703	12.770	ng	0.00
45) 2-Fluorobiphenyl	13.069	172	154335	9.455	ng	-0.01
79) Terphenyl-d14	19.774	244	245828	8.739	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	10.651	128	110346	5.647	ng	# 91
37) 2-Methylnaphthalene	12.269	142	735139	53.195	ng	98
38) 1-Methylnaphthalene	12.487	142	488679	36.328	ng	99
46) 1,1'-Biphenyl	13.281	154	34985	2.034	ng	99
52) Acenaphthene	14.516	154	49170	3.937	ng	93
55) Dibenzofuran	14.851	168	57459	2.692	ng	# 1
58) Fluorene	15.504	166	80506	4.642	ng	# 85
71) Phenanthrene	17.251	178	290775	11.042	ng	96

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

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