

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122921\
 Data File : BP008596.D
 Acq On : 29 Dec 2021 15:50
 Operator : CG/JU
 Sample : M5219-01DL 50X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 M21013-N48854-1040DL

Quant Time: Dec 29 16:20:34 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP122821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 29 13:29:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	569317	20.000	ng	0.00
21) Naphthalene-d8	10.704	136	2105896	20.000	ng	0.00
39) Acenaphthene-d10	14.534	164	1361235	20.000	ng	0.00
64) Phenanthrene-d10	17.281	188	2830108	20.000	ng	0.00
76) Chrysene-d12	21.357	240	2986639	20.000	ng	0.00
86) Perylene-d12	23.786	264	3263242	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	56605	1.815	ng	0.00
7) Phenol-d6	7.058	99	38475	0.886	ng	0.00
23) Nitrobenzene-d5	9.057	82	66158	1.869	ng	0.00
42) 2,4,6-Tribromophenol	16.016	330	39395	2.139	ng	0.00
45) 2-Fluorobiphenyl	13.157	172	215638	2.393	ng	0.00
79) Terphenyl-d14	19.833	244	316922	1.032	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	7.087	94	117694	2.685	ng	99
17) 2-Methylphenol	8.340	107	309077	10.134	ng	99
20) 3+4-Methylphenols	8.669	107	356211	8.313	ng	90
27) 2,4-Dimethylphenol	9.910	122	1087782	37.646	ng	# 90
38) 1-Methylnaphthalene	12.581	142	170848	2.124	ng	99

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

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