

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011521\
 Data File : BP004586.D
 Acq On : 15 Jan 2021 20:36
 Operator : CG/JU
 Sample : M1124-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ETGI-358

Quant Time: Jan 15 22:00:52 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP011221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 12 15:19:32 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	180264	20.00	ng	-0.01
21) Naphthalene-d8	10.598	136	702019	20.00	ng	#-0.01
39) Acenaphthene-d10	14.457	164	445607	20.00	ng	-0.01
64) Phenanthrene-d10	17.216	188	987585	20.00	ng	-0.01
76) Chrysene-d12	21.309	240	1025339	20.00	ng	-0.01
86) Perylene-d12	23.656	264	1146057	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.398	112	1446893	139.90	ng	0.00
7) Phenol-d6	6.987	99	1623625	114.62	ng	0.00
23) Nitrobenzene-d5	8.963	82	1088586	87.67	ng	-0.01
42) 2,4,6-Tribromophenol	15.957	330	1088592	140.51	ng	0.00
45) 2-Fluorobiphenyl	13.075	172	2993017	87.32	ng	0.00
79) Terphenyl-d14	19.786	244	4440400	79.12	ng	0.00
Target Compounds						
10) Phenol	7.010	94	141767	10.44	ng	94
50) Dimethylphthalate	13.910	163	243365	6.59	ng	99

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

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