

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112520\
 Data File : BP004106.D
 Acq On : 25 Nov 2020 20:12
 Operator : CG/JU
 Sample : L4858-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PHC-340EM-013

Quant Time: Nov 26 00:32:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 23 16:42:41 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.263	152	111228	20.00	ng	0.00
21) Naphthalene-d8	11.093	136	437631	20.00	ng	# 0.00
39) Acenaphthene-d10	14.886	164	275882	20.00	ng	0.00
64) Phenanthrene-d10	17.622	188	615788	20.00	ng	# 0.00
76) Chrysene-d12	21.657	240	617165	20.00	ng	0.00
86) Perylene-d12	24.127	264	697502	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.799	112	874075	131.16	ng	0.00
7) Phenol-d6	7.452	99	989521	103.43	ng	0.00
23) Nitrobenzene-d5	9.428	82	814644	91.17	ng	0.00
42) 2,4,6-Tribromophenol	16.375	330	524868	152.11	ng	0.00
45) 2-Fluorobiphenyl	13.510	172	1760972	89.24	ng	0.00
79) Terphenyl-d14	20.121	244	2345111	73.41	ng	0.00
Target Compounds						
10) Phenol	7.475	94	86898	9.41	ng	Qvalue 78

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

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