

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040221\
 Data File : BP005161.D
 Acq On : 02 Apr 2021 16:11
 Operator : CG/JU
 Sample : M1858-09
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 B4B

Quant Time: Apr 02 16:52:35 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 01 12:23:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	30225	20.00	ng	# 0.00
21) Naphthalene-d8	10.493	136	120338	20.00	ng	# 0.00
39) Acenaphthene-d10	14.357	164	72143	20.00	ng	0.00
64) Phenanthrene-d10	17.122	188	148177	20.00	ng	# 0.00
76) Chrysene-d12	21.216	240	141330	20.00	ng	-0.01
86) Perylene-d12	23.504	264	146611	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.305	112	317622	161.72	ng	0.00
7) Phenol-d6	6.887	99	413747	147.07	ng	0.00
23) Nitrobenzene-d5	8.863	82	268095	95.59	ng	0.00
42) 2,4,6-Tribromophenol	15.863	330	131444	139.84	ng	0.00
45) 2-Fluorobiphenyl	12.975	172	503823	94.47	ng	0.00
79) Terphenyl-d14	19.698	244	678132	87.35	ng	0.00
Target Compounds						
10) Phenol	6.911	94	6396	2.22	ng	Qvalue 81
32) Benzoic acid	9.769	122	9923	7.25	ng	# 70
50) Dimethylphthalate	13.816	163	21740	3.91	ng	97

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

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