

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092321\
 Data File : BP007148.D
 Acq On : 24 Sep 2021 04:24
 Operator : CG/JU
 Sample : M3881-01 10X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OK-01-092221

Quant Time: Sep 24 05:03:50 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 23 11:17:25 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	343461	20.000	ng	0.00
21) Naphthalene-d8	10.693	136	1360452	20.000	ng	# 0.00
39) Acenaphthene-d10	14.522	164	851669	20.000	ng	0.00
64) Phenanthrene-d10	17.275	188	1704320	20.000	ng	0.00
76) Chrysene-d12	21.357	240	1537581	20.000	ng	# 0.00
86) Perylene-d12	23.774	264	1699682	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	169352	8.254	ng	0.00
7) Phenol-d6	7.063	99	208789	7.445	ng	0.00
23) Nitrobenzene-d5	9.052	82	116398	4.983	ng	0.00
42) 2,4,6-Tribromophenol	16.010	330	89031	8.063	ng	0.00
45) 2-Fluorobiphenyl	13.145	172	333603	5.612	ng	0.00
79) Terphenyl-d14	19.827	244	454678	5.667	ng	0.00

Target Compounds	Qvalue

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

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