

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092321\
 Data File : BP007132.D
 Acq On : 23 Sep 2021 19:20
 Operator : CG/JU
 Sample : M3797-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 VB16193

Quant Time: Sep 24 01:01:30 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	298212	20.000	ng	0.00
21) Naphthalene-d8	10.693	136	1164030	20.000	ng	# 0.00
39) Acenaphthene-d10	14.522	164	752661	20.000	ng	0.00
64) Phenanthrene-d10	17.269	188	1543232	20.000	ng	# 0.00
76) Chrysene-d12	21.351	240	1585719	20.000	ng	# 0.00
86) Perylene-d12	23.762	264	1745167	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	1683628	94.507	ng	0.00
7) Phenol-d6	7.063	99	2040782	83.815	ng	0.00
23) Nitrobenzene-d5	9.051	82	1142522	57.160	ng	0.00
42) 2,4,6-Tribromophenol	16.010	330	974306	99.848	ng	0.00
45) 2-Fluorobiphenyl	13.145	172	1902395	36.210	ng	0.00
79) Terphenyl-d14	19.827	244	2853199	34.479	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.310	178	223207	2.697	ng	98
75) Fluoranthene	19.274	202	400414	4.200	ng	95
78) Pyrene	19.627	202	339212	3.260	ng	98

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

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