

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092321\
 Data File : BP007146.D
 Acq On : 24 Sep 2021 03:16
 Operator : CG/JU
 Sample : M3798-08DL 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 NAPL-06-(8-10)-END-POINTDL

Quant Time: Sep 24 04:10:50 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	7.899	152	410900	20.000	ng	0.00
21) Naphthalene-d8	10.693	136	1785449	20.000	ng	# 0.00
39) Acenaphthene-d10	14.522	164	1148580	20.000	ng	0.00
64) Phenanthrene-d10	17.275	188	2276826	20.000	ng	0.00
76) Chrysene-d12	21.357	240	2153314	20.000	ng	0.00
86) Perylene-d12	23.774	264	2368017	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	321746	13.107	ng	0.00
7) Phenol-d6	7.063	99	366720	10.931	ng	0.00
23) Nitrobenzene-d5	9.057	82	288676	9.416	ng	0.00
42) 2,4,6-Tribromophenol	16.010	330	176340	11.842	ng	0.00
45) 2-Fluorobiphenyl	13.145	172	607314	7.575	ng	0.00
79) Terphenyl-d14	19.827	244	826475	7.355	ng	0.00
Target Compounds						
31) Naphthalene	10.746	128	265132	2.910	ng	Qvalue 99
84) Bis(2-ethylhexyl)phtha...	21.269	149	453113	5.571	ng	97

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

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