

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081122\
 Data File : BM036218.D
 Acq On : 11 Aug 2022 20:36
 Operator : CG/JU
 Sample : N4102-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PPE-DRUMS-0808

Quant Time: Aug 12 02:18:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 10 22:19:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	320941	20.000	ng	0.00
21) Naphthalene-d8	10.633	136	1382230	20.000	ng	0.00
39) Acenaphthene-d10	14.469	164	844258	20.000	ng	0.00
64) Phenanthrene-d10	17.210	188	1763256	20.000	ng	0.00
76) Chrysene-d12	21.392	240	1986184	20.000	ng	0.00
86) Perylene-d12	23.738	264	1573110	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	3155029	159.379	ng	0.00
7) Phenol-d6	7.034	99	3771825	149.743	ng	0.00
23) Nitrobenzene-d5	9.004	82	2616810	105.974	ng	0.00
42) 2,4,6-Tribromophenol	15.963	330	1886758	190.500	ng	0.00
45) 2-Fluorobiphenyl	13.098	172	5781554	101.110	ng	0.00
79) Terphenyl-d14	19.839	244	8809225	87.483	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.681	128	4788433	69.679	ng	99
37) 2-Methylnaphthalene	12.292	142	1166609	25.511	ng	97
38) 1-Methylnaphthalene	12.516	142	948306	21.175	ng	98
46) 1,1'-Biphenyl	13.304	154	153380	2.493	ng	99
52) Acenaphthene	14.533	154	410443	8.620	ng	100
58) Fluorene	15.515	166	241182	4.288	ng	97
71) Phenanthrene	17.251	178	570510	6.351	ng	100

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

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