

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071422\
 Data File : BM035810.D
 Acq On : 14 Jul 2022 13:47
 Operator : CG/JU
 Sample : N3651-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SED-1-(6-12)

Quant Time: Jul 15 01:09:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM071322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 14 01:21:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	319848	20.000	ng	0.00
21) Naphthalene-d8	10.716	136	1318342	20.000	ng	0.00
39) Acenaphthene-d10	14.539	164	757212	20.000	ng	0.00
64) Phenanthrene-d10	17.280	188	1468634	20.000	ng	0.00
76) Chrysene-d12	21.456	240	1260860	20.000	ng	0.00
86) Perylene-d12	23.839	264	1259186	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	2075119	105.381	ng	0.00
7) Phenol-d6	7.075	99	2558430	103.869	ng	0.00
23) Nitrobenzene-d5	9.075	82	1534134	66.525	ng	0.00
42) 2,4,6-Tribromophenol	16.027	330	863635	112.313	ng	0.00
45) 2-Fluorobiphenyl	13.175	172	3144607	57.917	ng	0.00
79) Terphenyl-d14	19.903	244	4020593	61.900	ng	0.00
Target Compounds						
74) Di-n-butylphthalate	18.257	149	3827	2.309	ng	# 87
84) Bis(2-ethylhexyl)phtha...	21.386	149	6674	3.382	ng	# 92

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

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