

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080522\
 Data File : BM036132.D
 Acq On : 05 Aug 2022 23:04
 Operator : CG/JU
 Sample : N3961-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RVE-173

Quant Time: Aug 06 01:14:12 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 03:36:25 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	269680	20.00	ng	0.00
21) Naphthalene-d8	10.657	136	1089708	20.00	ng	0.00
39) Acenaphthene-d10	14.492	164	622531	20.00	ng	0.00
64) Phenanthrene-d10	17.239	188	1231044	20.00	ng	0.00
76) Chrysene-d12	21.415	240	1135675	20.00	ng	0.00
86) Perylene-d12	23.774	264	1199980	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.445	112	1804845	108.50	ng	0.02
7) Phenol-d6	7.069	99	2228967	105.31	ng	0.04
23) Nitrobenzene-d5	9.028	82	1348480	69.27	ng	0.00
42) 2,4,6-Tribromophenol	15.986	330	851567	116.60	ng	0.00
45) 2-Fluorobiphenyl	13.121	172	3038792	72.07	ng	0.00
79) Terphenyl-d14	19.862	244	4274229	74.23	ng	0.00
Target Compounds						
74) Di-n-butylphthalate	18.209	149	443005	6.03	ng	Qvalue 99
75) Fluoranthene	19.292	202	254201	3.62	ng	97
78) Pyrene	19.656	202	291419	4.14	ng	99
88) Benzo(b)fluoranthene	23.056	252	234405	3.33	ng	# 90
90) Benzo(a)pyrene	23.668	252	161127	2.73	ng	# 91

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

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