

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080122\
 Data File : BF129471.D
 Acq On : 01 Aug 2022 13:55
 Operator : CG\JU
 Sample : N3964-07DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HMGP-WATER-1DL

Quant Time: Aug 02 01:15:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 01:11:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	177337	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	676808	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	313745	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	438856	20.000	ng	0.00
76) Chrysene-d12	14.045	240	374617	20.000	ng	0.00
86) Perylene-d12	15.527	264	449131	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	130577	11.995	ng	0.01
7) Phenol-d6	6.528	99	96032	7.018	ng	0.00
23) Nitrobenzene-d5	7.434	82	207747	16.756	ng	-0.01
42) 2,4,6-Tribromophenol	10.704	330	57897	19.194	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	375964	18.537	ng	0.00
79) Terphenyl-d14	12.980	244	301585	15.188	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.181	128	965595	28.081	ng	99
37) 2-Methylnaphthalene	8.869	142	135739	6.074	ng	98
38) 1-Methylnaphthalene	8.969	142	145590	6.749	ng	98
52) Acenaphthene	9.945	154	72483	3.946	ng	97
58) Fluorene	10.457	166	69013	3.406	ng	97
71) Phenanthrene	11.428	178	177935	7.428	ng	98
75) Fluoranthene	12.616	202	49862	2.054	ng	98
78) Pyrene	12.845	202	64037	2.044	ng	96

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

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