

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF032922\
 Data File : BF127564.D
 Acq On : 29 Mar 2022 14:38
 Operator : CG\JU
 Sample : N2154-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-4

Quant Time: Mar 29 15:56:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF032122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 29 07:13:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	82791	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	299199	20.000	ng	0.00
39) Acenaphthene-d10	9.869	164	163573	20.000	ng	0.00
64) Phenanthrene-d10	11.363	188	259399	20.000	ng	0.00
76) Chrysene-d12	14.015	240	163373	20.000	ng	0.00
86) Perylene-d12	15.498	264	185212	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	311545	60.066	ng	0.00
7) Phenol-d6	6.475	99	258522	36.464	ng	0.00
23) Nitrobenzene-d5	7.404	82	582339	83.197	ng	0.00
42) 2,4,6-Tribromophenol	10.669	330	196754	112.917	ng	0.00
45) 2-Fluorobiphenyl	9.192	172	943240	84.204	ng	0.00
79) Terphenyl-d14	12.945	244	706207	70.763	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.492	94	31184	4.039	ng	91
32) Benzoic acid	7.969	122	169615	67.300	ng	93
36) 4-Chloro-3-methylphenol	8.686	107	37686	7.265	ng	92

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

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