

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF032522\
 Data File : BF127527.D
 Acq On : 25 Mar 2022 15:50
 Operator : CG\JU
 Sample : N2076-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-20220322

Quant Time: Mar 25 17:27:03 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 22 02:16:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.845	152	88911	20.000	ng	0.00
21) Naphthalene-d8	8.133	136	316306	20.000	ng	0.00
39) Acenaphthene-d10	9.886	164	194325	20.000	ng	0.00
64) Phenanthrene-d10	11.380	188	363295	20.000	ng	0.00
76) Chrysene-d12	14.027	240	292256	20.000	ng	0.00
86) Perylene-d12	15.509	264	204371	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	633780	113.783	ng	0.01
7) Phenol-d6	6.486	99	812056	106.655	ng	0.00
23) Nitrobenzene-d5	7.416	82	640388	86.542	ng	0.00
42) 2,4,6-Tribromophenol	10.686	330	272127	131.459	ng	0.00
45) 2-Fluorobiphenyl	9.204	172	1085139	81.541	ng	0.00
79) Terphenyl-d14	12.968	244	1280231	71.710	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.498	94	39655	4.782	ng	86
27) 2,4-Dimethylphenol	7.798	122	10623	2.386	ng	94
31) Naphthalene	8.169	128	4084638	243.232	ng	97
37) 2-Methylnaphthalene	8.839	142	368084	34.088	ng	99
38) 1-Methylnaphthalene	8.939	142	322920	30.243	ng	98
46) 1,1'-Biphenyl	9.304	154	37682	2.567	ng	99
50) Dimethylphthalate	9.598	163	52881	3.912	ng	96
52) Acenaphthene	9.916	154	67288	6.601	ng	98
71) Phenanthrene	11.404	178	76013	4.013	ng	97

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

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