

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110922\
 Data File : BF131100.D
 Acq On : 09 Nov 2022 22:06
 Operator : CG\JU
 Sample : N5340-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-2

Quant Time: Nov 10 04:33:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 10 04:15:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.893	152	91309	20.000	ng	0.00
21) Naphthalene-d8	8.181	136	324821	20.000	ng	0.00
39) Acenaphthene-d10	9.939	164	168703	20.000	ng	0.00
64) Phenanthrene-d10	11.433	188	271606	20.000	ng	0.00
76) Chrysene-d12	14.074	240	182401	20.000	ng	# 0.00
86) Perylene-d12	15.568	264	190623	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	272552	50.762	ng	0.01
7) Phenol-d6	6.522	99	164457	23.429	ng	0.00
23) Nitrobenzene-d5	7.463	82	674026	93.625	ng	0.00
42) 2,4,6-Tribromophenol	10.739	330	234032	134.231	ng	0.00
45) 2-Fluorobiphenyl	9.257	172	1086249	89.884	ng	0.00
79) Terphenyl-d14	13.022	244	921659	88.594	ng	0.00
Target Compounds						
						Qvalue
4) n-Nitrosodimethylamine	3.416	42	49075	11.723	ng	# 93
10) Phenol	6.540	94	107019	13.042	ng	94
15) Benzyl Alcohol	7.051	79	521547	97.146	ng	98
22) Acetophenone	7.298	105	141077	17.355	ng	# 96
26) 2-Nitrophenol	7.792	139	23881	7.718	ng	94
32) Benzoic acid	8.198	122	1860373	497.982	ng	94
36) 4-Chloro-3-methylphenol	8.739	107	91868	16.364	ng	# 89
37) 2-Methylnaphthalene	8.892	142	51470	4.499	ng	98
38) 1-Methylnaphthalene	8.992	142	47340	4.225	ng	100
52) Acenaphthene	9.969	154	24962	2.587	ng	99
71) Phenanthrene	11.457	178	42387	2.838	ng	# 92

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

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