

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041422\
 Data File : BF127795.D
 Acq On : 14 Apr 2022 15:49
 Operator : CG\JU
 Sample : N2389-07
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 33637-8

Quant Time: Apr 15 05:06:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 16:40:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.957	152	102009	20.000	ng	0.00
21) Naphthalene-d8	8.239	136	333906	20.000	ng	# 0.00
39) Acenaphthene-d10	10.004	164	141070	20.000	ng	0.00
64) Phenanthrene-d10	11.492	188	260510	20.000	ng	# 0.00
76) Chrysene-d12	14.139	240	313388	20.000	ng	0.00
86) Perylene-d12	15.662	264	341921	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.592	112	623503	99.318	ng	0.01
7) Phenol-d6	6.592	99	662642	80.684	ng	0.00
23) Nitrobenzene-d5	7.522	82	559670	74.724	ng	0.00
42) 2,4,6-Tribromophenol	10.804	330	184085	122.626	ng	0.00
45) 2-Fluorobiphenyl	9.316	172	897415	89.994	ng	0.00
79) Terphenyl-d14	13.080	244	1054928	55.563	ng	0.00
Target Compounds						
						Qvalue
9) Benzaldehyde	6.510	77	60929	8.776	ng	95
10) Phenol	6.604	94	110558	13.463	ng	94
15) Benzyl Alcohol	7.104	79	65395	9.108	ng	98
17) 2-Methylphenol	7.204	107	35008	6.320	ng	99
22) Acetophenone	7.363	105	198157	22.085	ng	# 81
27) 2,4-Dimethylphenol	7.892	122	14306	2.800	ng	# 88
31) Naphthalene	8.263	128	1068302	61.057	ng	98
32) Benzoic acid	8.157	122	1726344	465.789	ng	98
37) 2-Methylnaphthalene	8.951	142	27008	2.308	ng	99

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

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