

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061521\
 Data File : BF124334.D
 Acq On : 15 Jun 2021 21:24
 Operator : JU/CG
 Sample : M2674-11DL 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TMR-DS-05-060921DL

Quant Time: Jun 16 01:28:08 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 15 20:09:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	31990	20.00	ng	0.00
21) Naphthalene-d8	8.069	136	126753	20.00	ng	0.00
39) Acenaphthene-d10	9.827	164	76318	20.00	ng	0.00
64) Phenanthrene-d10	11.316	188	143251	20.00	ng	# 0.00
76) Chrysene-d12	13.957	240	100759	20.00	ng	0.00
86) Perylene-d12	15.415	264	89006	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	48105	22.64	ng	0.00
7) Phenol-d6	6.439	99	52478	18.37	ng	0.00
23) Nitrobenzene-d5	7.351	82	58956	18.45	ng	0.00
42) 2,4,6-Tribromophenol	10.616	330	26121	24.16	ng	0.00
45) 2-Fluorobiphenyl	9.145	172	99565	16.42	ng	0.00
79) Terphenyl-d14	12.898	244	117307	15.98	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.457	94	35418	11.47	ng	97
17) 2-Methylphenol	7.057	107	60982	31.72	ng	91
20) 3+4-Methylphenols	7.216	107	114222	44.85	ng	# 71
22) Acetophenone	7.192	105	9442	2.61	ng	# 94
27) 2,4-Dimethylphenol	7.739	122	107895	53.43	ng	91
31) Naphthalene	8.092	128	16452	2.17	ng	# 95

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

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