

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061521\
 Data File : BF124336.D
 Acq On : 15 Jun 2021 22:30
 Operator : JU/CG
 Sample : M2674-12DL 10X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TMR-DS-06-060921DL

Quant Time: Jun 16 01:28:29 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	33966	20.00	ng	0.00
21) Naphthalene-d8	8.069	136	126364	20.00	ng	# 0.00
39) Acenaphthene-d10	9.822	164	76863	20.00	ng	0.00
64) Phenanthrene-d10	11.310	188	138206	20.00	ng	0.00
76) Chrysene-d12	13.957	240	87399	20.00	ng	0.00
86) Perylene-d12	15.409	264	82790	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	24173	10.71	ng	0.00
7) Phenol-d6	6.439	99	26696	8.80	ng	0.00
23) Nitrobenzene-d5	7.345	82	30376	9.53	ng	-0.01
42) 2,4,6-Tribromophenol	10.616	330	12992	11.93	ng	0.00
45) 2-Fluorobiphenyl	9.145	172	48600	7.96	ng	0.00
79) Terphenyl-d14	12.892	244	51247	8.05	ng	-0.01
Target Compounds						
10) Phenol	6.451	94	30120	9.19	ng	Qvalue 93
17) 2-Methylphenol	7.051	107	40408	19.80	ng	90
20) 3+4-Methylphenols	7.216	107	100038	36.99	ng	# 69
27) 2,4-Dimethylphenol	7.733	122	69619	34.58	ng	92

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

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