

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061521\
 Data File : BF124332.D
 Acq On : 15 Jun 2021 20:19
 Operator : JU/CG
 Sample : M2674-09 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TMR-DS-03-060921

Quant Time: Jun 15 20:40:09 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	35147	20.00	ng	0.00
21) Naphthalene-d8	8.069	136	129538	20.00	ng	# 0.00
39) Acenaphthene-d10	9.822	164	73111	20.00	ng	0.00
64) Phenanthrene-d10	11.310	188	129467	20.00	ng	0.00
76) Chrysene-d12	13.957	240	81385	20.00	ng	0.00
86) Perylene-d12	15.410	264	72794	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	48925	20.95	ng	0.00
7) Phenol-d6	6.440	99	56952	18.15	ng	0.00
23) Nitrobenzene-d5	7.345	82	55216	16.91	ng	-0.01
42) 2,4,6-Tribromophenol	10.616	330	22764	21.98	ng	0.00
45) 2-Fluorobiphenyl	9.145	172	101637	17.49	ng	0.00
79) Terphenyl-d14	12.898	244	97813	16.50	ng	0.00
Target Compounds						
10) Phenol	6.457	94	37192	10.97	ng	Qvalue 93
17) 2-Methylphenol	7.057	107	41280	19.55	ng	97
20) 3+4-Methylphenols	7.216	107	104416	37.31	ng	# 69
27) 2,4-Dimethylphenol	7.734	122	66798	32.37	ng	94

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

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