

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111920\
 Data File : BF122389.D
 Acq On : 19 Nov 2020 19:39
 Operator : JU/CG
 Sample : L4779-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1623

Manual Integrations
 APPROVED

mohammad
 11/24/2020 8:54:22 AM

Quant Time: Nov 20 00:03:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 10 12:25:42 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	0.000	152	0	0.00	ng	-6.87
21) Naphthalene-d8	8.181	136	32542m	20.00	ng	0.03
39) Acenaphthene-d10	9.945	164	2687	20.00	ng	# 0.04
64) Phenanthrene-d10	11.445	188	1892	20.00	ng	# 0.05
76) Chrysene-d12	14.110	240	717	20.00	ng	# 0.08
86) Perylene-d12	0.000	264	0m	20.00	ng	-15.50
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	1303805	0.00	ng	0.00
7) Phenol-d6	6.487	99	1232797	0.00	ng	0.00
23) Nitrobenzene-d5	7.428	82	1407538	2647.93	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	801932	27807.15	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	2554950	14856.66	ng	0.00
79) Terphenyl-d14	12.980	244	3111360	91930.64	ng	0.00
Target Compounds						Qvalue
22) Acetophenone	7.269	105	10371	12.24	ng	# 1
25) Isophorone	7.698	82	64606	54.52	ng	# 72
27) 2,4-Dimethylphenol	7.804	122	9271	16.59	ng	82
28) bis(2-Chloroethoxy)met...	7.898	93	47758	65.89	ng	# 1
32) Benzoic acid	8.028	122	90323m	251.46	ng	
46) 1,1'-Biphenyl	9.328	154	5820	27.21	ng	95
49) Acenaphthylene	9.822	152	1062	4.07	ng	# 33
50) Dimethylphthalate	9.622	163	331173	1733.94	ng	99
70) Pentachlorophenol	11.198	266	3161	228.40	ng	# 88
78) Pyrene	12.827	202	2433m	48.30	ng	

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

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