

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110218\
 Data File : BF110430.D
 Acq On : 2 Nov 2018 23:42
 Operator : JU/SJ
 Sample : J5769-05 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WAGARAW-RD-3

Manual Integrations
 APPROVED

Sohil
 11/5/2018 10:32:51 AM

Quant Time: Nov 03 01:05:00 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 16:06:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	92034	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	355061	20.00	ng	0.00
39) Acenaphthene-d10	10.00	164	140384	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	221429	20.00	ng	0.00
76) Chrysene-d12	14.13	240	154152	20.00	ng	0.00
87) Perylene-d12	15.66	264	112012	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	265099	46.91	ng	0.00
7) Phenol-d6	6.57	99	337313	46.66	ng	0.00
23) Nitrobenzene-d5	7.52	82	195029	32.58	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	55880	40.18	ng	0.00
45) 2-Fluorobiphenyl	9.31	172	372869	41.71	ng	0.00
79) Terphenyl-d14	13.07	244	276964	37.37	ng	0.00
Target Compounds						
10) Phenol	6.59	94	20623	2.669	ng	Qvalue # 77
50) Dimethylphthalate	9.70	163	49271	4.755	ng	99
71) Phenanthrene	11.51	178	41144	3.551	ng	97
75) Fluoranthene	12.70	202	66379	5.645	ng	99
78) Pyrene	12.93	202	69719	6.309	ng	98
81) Benzo(a)anthracene	14.12	228	30282	3.313	ng	# 91
83) Chrysene	14.16	228	32218	3.711	ng	96
88) Benzo(b)fluoranthene	15.20	252	38479m	5.949	ng	
90) Benzo(a)pyrene	15.59	252	21654	3.638	ng	# 93
92) Benzo(g,h,i)perylene	17.70	276	16274	3.216	ng	# 90

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

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