

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110218\
 Data File : BF110434.D
 Acq On : 3 Nov 2018 1:32
 Operator : JU/SJ
 Sample : J5769-17
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GOFFLE-RD-2

Manual Integrations
 APPROVED

Sohil
 11/5/2018 10:32:56 AM

Quant Time: Nov 03 03:38:59 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	81499	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	299200	20.00	ng	0.00
39) Acenaphthene-d10	10.00	164	122355	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	206840	20.00	ng	0.00
76) Chrysene-d12	14.13	240	136696	20.00	ng	0.00
87) Perylene-d12	15.66	264	100498	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	513816	102.67	ng	0.00
7) Phenol-d6	6.58	99	635360	99.25	ng	0.00
23) Nitrobenzene-d5	7.52	82	377169	74.77	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	114263	94.28	ng	0.00
45) 2-Fluorobiphenyl	9.31	172	656452	84.25	ng	0.00
79) Terphenyl-d14	13.07	244	509906	77.58	ng	0.00
Target Compounds						
10) Phenol	6.59	94	37249	5.444	ng	Qvalue # 68
50) Dimethylphthalate	9.70	163	87049	9.638	ng	99
71) Phenanthrene	11.52	178	38685	3.574	ng	97
75) Fluoranthene	12.70	202	95391	8.685	ng	100
78) Pyrene	12.94	202	91545	9.341	ng	98
81) Benzo(a)anthracene	14.12	228	39502m	4.873	ng	
83) Chrysene	14.16	228	37783	4.907	ng	97
86) Indeno(1,2,3-cd)pyrene	17.23	276	14565	2.280	ng	# 83
88) Benzo(b)fluoranthene	15.20	252	40528m	6.983	ng	
89) Benzo(k)fluoranthene	15.23	252	12072m	2.116	ng	
90) Benzo(a)pyrene	15.60	252	26725	5.005	ng	# 96
92) Benzo(a,h,i)perylene	17.70	276	16621	3.661	ng	# 88

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

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