

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112818\
 Data File : BF111059.D
 Acq On : 28 Nov 2018 21:34
 Operator : JU/SJ
 Sample : J6079-04
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 J-4

Manual Integrations
 APPROVED

mohammad
 11/29/2018 1:58:37 PM

Quant Time: Nov 29 03:29:24 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 27 15:32:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	54379	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	216337	20.00	ng	0.00
39) Acenaphthene-d10	9.97	164	86795	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	126707	20.00	ng	0.00
76) Chrysene-d12	14.13	240	109889	20.00	ng	0.00
87) Perylene-d12	15.66	264	85628	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	309322	94.33	ng	0.00
7) Phenol-d6	6.56	99	400917	93.82	ng	0.00
23) Nitrobenzene-d5	7.50	82	251620	66.56	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	68601	79.58	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	425701	71.06	ng	0.00
79) Terphenyl-d14	13.05	244	272974	48.36	ng	0.00
Target Compounds						
10) Phenol	6.57	94	19547	4.062	ng	83
50) Dimethylphthalate	9.69	163	33138	5.206	ng	98
71) Phenanthrene	11.50	178	26859	4.002	ng	98
75) Fluoranthene	12.70	202	36038	5.225	ng	97
78) Pyrene	12.93	202	40987	5.107	ng	97
81) Benzo(a)anthracene	14.12	228	15801m	2.436	ng	
83) Chrysene	14.16	228	19531	3.058	ng	97
84) Bis(2-ethylhexyl)phthalate	14.06	149	13742	2.846	ng	96
88) Benzo(b)fluoranthene	15.22	252	16495m	3.279	ng	
90) Benzo(a)pyrene	15.60	252	11246	2.402	ng	94

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

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