

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061518\
 Data File : BF106491.D
 Acq On : 15 Jun 2018 15:01
 Operator : JU/SJ
 Sample : J3471-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 #1

Manual Integrations
 APPROVED

Sohil
 6/18/2018 10:38:22 AM

Quant Time: Jun 15 23:57:44 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 15 11:11:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	113196	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	422463	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	181922	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	294738	20.00	ng	0.00
75) Chrysene-d12	14.15	240	301304	20.00	ng	0.00
86) Perylene-d12	15.69	264	275949	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	678949	101.52	ng	0.00
7) Phenol-d6	6.61	99	823665	97.48	ng	0.00
23) Nitrobenzene-d5	7.52	82	552170	76.89	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	178462	101.95	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	840598	80.76	ng	0.00
78) Terphenyl-d14	13.08	244	687586	56.68	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	99503	7.656	ng	98
70) Phenanthrene	11.52	178	39961	2.768	ng	98
74) Fluoranthene	12.71	202	95799	6.223	ng	96
77) Pyrene	12.94	202	92396	4.898	ng	98
80) Benzo(a)anthracene	14.14	228	53807	3.001	ng	97
82) Chrysene	14.18	228	57162	3.492	ng	98
83) Bis(2-ethylhexyl)phthalate	14.11	149	80429	7.944	ng	100
87) Benzo(b)fluoranthene	15.24	252	81310m	4.874	ng	
89) Benzo(a)pyrene	15.62	252	48585	3.238	ng	# 92
91) Benzo(g,h,i)perylene	17.74	276	25513	2.099	ng	91

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

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