

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083118\
 Data File : BF108698.D
 Acq On : 31 Aug 2018 19:23
 Operator : JU/SJ
 Sample : J4720-05 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BZTRANS-5

Manual Integrations
 APPROVED

Sohil
 9/4/2018 3:53:24 PM

Quant Time: Sep 01 04:08:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 30 15:06:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	65677	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	313148	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	134286	20.00	ng	0.00
63) Phenanthrene-d10	11.54	188	242230	20.00	ng	0.00
75) Chrysene-d12	14.19	240	232379	20.00	ng	0.00
86) Perylene-d12	15.75	264	173678	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.85	112	25209m	6.67	ng	0.20
7) Phenol-d6	6.70	99	61942m	11.44	ng	0.05
23) Nitrobenzene-d5	7.62	82	52438m	8.50	ng	0.05
41) 2,4,6-Tribromophenol	10.84	330	17206	9.71	ng	0.00
44) 2-Fluorobiphenyl	9.35	172	107135	10.73	ng	0.00
78) Terphenyl-d14	13.12	244	108114	9.04	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.75	163	32725	3.095	ng	Qvalue # 99
70) Phenanthrene	11.57	178	68845	5.631	ng	97
74) Fluoranthene	12.76	202	90769	6.595	ng	94
77) Pyrene	12.99	202	70260	4.022	ng	98
82) Chrysene	14.22	228	31748	2.330	ng	98
87) Benzo(b)fluoranthene	15.29	252	34889m	3.270	ng	
89) Benzo(a)pyrene	15.69	252	21692	2.250	ng	# 93

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

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