

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF021218\
 Data File : BF102920.D
 Acq On : 12 Feb 2018 15:07
 Operator : SJ/JU
 Sample : J1491-01RX
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1-1-4-DENSE-GRADE-AGG-RCARX

Manual Integrations
 APPROVED

Sohil
 2/13/2018 11:25:35 AM

Quant Time: Feb 13 00:39:25 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 12:44:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	192975	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	784526	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	310801	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	442903	20.00	ng	0.00
75) Chrysene-d12	14.06	240	361975	20.00	ng	0.00
86) Perylene-d12	15.54	264	324259	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	81855	6.44	ng	0.01
7) Phenol-d6	6.51	99	642770	42.01	ng	0.00
23) Nitrobenzene-d5	7.45	82	1094589	96.79	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	15409	6.16	ng	0.00
44) 2-Fluorobiphenyl	9.25	172	1529420	81.66	ng	0.00
78) Terphenyl-d14	13.00	244	1055486	69.04	ng	0.00
Target Compounds						
10) Phenol	6.52	94	129520	7.899	ng	95
49) Dimethylphthalate	9.63	163	146025	6.026	ng	99
70) Phenanthrene	11.43	178	93841	3.804	ng	97
74) Fluoranthene	12.63	202	101356	4.084	ng	98
77) Pyrene	12.86	202	93524	3.295	ng	99
80) Benzo(a)anthracene	14.05	228	47484m	2.086	ng	
87) Benzo(b)fluoranthene	15.10	252	52060m	2.476	ng	

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

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