

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082720\
 Data File : BN011669.D
 Acq On : 27 Aug 2020 18:06
 Operator : CG/JU
 Sample : L3737-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 KY038PS038-200818

Quant Time: Aug 28 00:45:13 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 27 14:34:48 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	1902	0.40	ng	0.00
7) Naphthalene-d8	10.53	136	7096	0.40	ng	0.00
13) Acenaphthene-d10	14.38	164	4177	0.40	ng	0.00
19) Phenanthrene-d10	17.13	188	9392	0.40	ng	0.00
29) Chrysene-d12	21.32	240	9586	0.40	ng	0.00
36) Perylene-d12	23.58	264	10167	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.34	112	922	0.17	ng	0.00
5) Phenol-d6	6.90	99	664	0.09	ng	0.00
8) Nitrobenzene-d5	8.88	82	3486	0.51	ng	0.00
11) 2-Methylnaphthalene-d10	12.12	152	3636	0.28	ng	0.00
14) 2,4,6-Tribromophenol	15.88	330	764	0.38	ng	0.00
15) 2-Fluorobiphenyl	13.00	172	9227	0.52	ng	0.00
27) Fluoranthene-d10	19.16	212	10101	0.34	ng	0.00
31) Terphenyl-d14	19.77	244	17990	0.76	ng	0.00
Target Compounds						
30) Pyrene	19.55	202	755	0.023	ng	# 93
34) Bis(2-ethylhexyl)phthalate	21.25	149	462	0.033	ng	# 99
37) Benzo(b)fluoranthene	22.90	252	800	0.021	ng	# 8

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

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