

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025167.D
 Acq On : 29 Feb 2020 14:30
 Operator : CG/JU
 Sample : L1507-14DL 5X
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD12DL

Manual Integrations
 APPROVED

Sohil
 3/3/2020 8:39:41 AM

Quant Time: Mar 02 01:58:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 02 01:15:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	2845	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	11256	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	7480	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	17937	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	19911	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	23026	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	1182	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	2635	0.04	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.50	128	112025	3.269	ng/ul#	82
5) 2-Methylnaphthalene	12.12	142	86990	3.518	ng/ul	99
7) Acenaphthylene	14.03	152	78387	1.818	ng/ul#	93
8) Acenaphthene	14.37	153	155634	5.371	ng/ul	98
9) Fluorene	15.36	166	133763	3.679	ng/ul#	95
12) Phenanthrene	17.09	178	705880	11.188	ng/ul	95
13) Anthracene	17.18	178	300591	4.916	ng/ul	95
15) Fluoranthene	19.11	202	2756925	30.083	ng/ul	96
17) Pyrene	19.48	202	5725372	63.815	ng/ul	97
18) Benzo(a)anthracene	21.22	228	1077115	12.530	ng/ul	98
19) Chrysene	21.27	228	1404012	16.397	ng/ul	98
21) Benzo(b)fluoranthene	22.79	252	1682744m	18.643	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	568541m	6.358	ng/ul	
23) Benzo(a)pyrene	23.34	252	677891	8.226	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.62	276	350908	3.399	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.62	278	93189	1.074	ng/ul	96
26) Benzo(a,h,i)perylene	26.28	276	221610	2.547	ng/ul#	92

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

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