

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022520\
 Data File : BM025077.D
 Acq On : 26 Feb 2020 13:15
 Operator : CG/JU
 Sample : L1541-02
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBCX7

Quant Time: Feb 26 13:49:26 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 : Tue Feb 25 11:03:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	2137	0.40	ng/ul	0.00
2) Naphthalene-d8	10.47	136	7957	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.32	164	4869	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	11337	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	11916	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	13228	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	5372	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	14884	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.52	128	7877	0.325	ng/ul#	84
5) 2-Methylnaphthalene	12.13	142	9991	0.572	ng/ul	100
7) Acenaphthylene	14.04	152	6759	0.241	ng/ul	96
9) Fluorene	15.37	166	11976	0.506	ng/ul	97
11) Pentachlorophenol	16.72	266	3109	0.898	ng/ul	97
12) Phenanthrene	17.11	178	13532	0.339	ng/ul	97
13) Anthracene	17.20	178	18996	0.492	ng/ul	96
15) Fluoranthene	19.12	202	14675	0.253	ng/ul	97
17) Pyrene	19.48	202	18789	0.350	ng/ul	96
18) Benzo(a)anthracene	21.23	228	24668	0.479	ng/ul	98
19) Chrysene	21.28	228	24430	0.477	ng/ul	98
21) Benzo(b)fluoranthene	22.79	252	39946	0.770	ng/ul	91
22) Benzo(k)fluoranthene	22.84	252	17937	0.349	ng/ul#	97
23) Benzo(a)pyrene	23.35	252	16872	0.356	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.66	278	40809	0.819	ng/ul#	94
26) Benzo(a,h,i)perylene	26.31	276	17646	0.353	ng/ul#	92

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

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