

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060820\
 Data File : BM026299.D
 Acq On : 08 Jun 2020 22:34
 Operator : JU/CG
 Sample : L2815-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OR-02-060520

Manual Integrations
 APPROVED

mohammad
 6/9/2020 12:21:58 PM

Quant Time: Jun 09 02:28:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 08 15:00:18 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	239734	20.00	ng	0.01
21) Naphthalene-d8	10.21	136	962678	20.00	ng	0.00
39) Acenaphthene-d10	14.10	164	531465	20.00	ng	0.00
64) Phenanthrene-d10	16.85	188	917271	20.00	ng	0.00
76) Chrysene-d12	21.06	240	736221	20.00	ng	0.00
86) Perylene-d12	23.17	264	796631	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.09	112	1465065	108.05	ng	0.01
7) Phenol-d6	6.66	99	2096416	106.77	ng	0.02
23) Nitrobenzene-d5	8.60	82	1391407	70.97	ng	0.00
42) 2,4,6-Tribromophenol	15.60	330	641937	99.71	ng	0.01
45) 2-Fluorobiphenyl	12.71	172	2863713	81.79	ng	0.00
79) Terphenyl-d14	19.50	244	2937836	77.46	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.57	163	245584	6.318	ng	99
71) Phenanthrene	16.89	178	708717	14.837	ng	98
72) Anthracene	16.99	178	176038	3.693	ng	98
73) Carbazole	17.26	167	90953	2.012	ng	98
75) Fluoranthene	18.92	202	1196876	21.855	ng	99
78) Pyrene	19.29	202	1043291	21.447	ng	99
81) Benzo(a)anthracene	21.04	228	476591	10.785	ng	98
83) Chrysene	21.09	228	448331	10.646	ng	96
87) Indeno(1,2,3-cd)pyrene	25.23	276	309725	6.721	ng	95
88) Benzo(b)fluoranthene	22.55	252	591828	12.350	ng	# 91
89) Benzo(k)fluoranthene	22.59	252	198209m	4.257	ng	
90) Benzo(a)pyrene	23.07	252	447002	10.502	ng	# 92
92) Benzo(a,h,i)perylene	25.87	276	324493	8.864	ng	98

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

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