

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012920\
 Data File : BM024653.D
 Acq On : 29 Jan 2020 16:53
 Operator : CG/JU
 Sample : L1278-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TONNELE-AVE-01

Manual Integrations
 APPROVED

mohammad
 1/30/2020 10:19:01 AM

Quant Time: Jan 30 03:46:40 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM012320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 23 14:57:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	401264	20.00	ng	0.00
21) Naphthalene-d8	10.20	136	1471635	20.00	ng	0.00
39) Acenaphthene-d10	14.09	164	856002	20.00	ng	0.00
64) Phenanthrene-d10	16.84	188	1497500	20.00	ng	0.00
76) Chrysene-d12	21.05	240	1233808	20.00	ng	0.00
87) Perylene-d12	23.16	264	1505266	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.09	112	2264742	107.34	ng	0.00
7) Phenol-d6	6.65	99	2831352	97.41	ng	0.00
23) Nitrobenzene-d5	8.58	82	1822036	60.72	ng	0.00
42) 2,4,6-Tribromophenol	15.59	330	1224882	87.97	ng	0.00
45) 2-Fluorobiphenyl	12.70	172	4622562	73.42	ng	0.00
79) Terphenyl-d14	19.50	244	4901647	74.17	ng	0.00
Target Compounds						
10) Phenol	6.67	94	144158	4.975	ng	98
50) Dimethylphthalate	13.56	163	601052	8.683	ng	100
71) Phenanthrene	16.88	178	410565	5.063	ng	98
75) Fluoranthene	18.91	202	706829	6.984	ng	98
78) Pyrene	19.27	202	670841	8.248	ng	99
81) Benzo(a)anthracene	21.03	228	379054	4.426	ng	96
83) Chrysene	21.09	228	366972	4.491	ng	97
86) Indeno(1,2,3-cd)pyrene	25.23	276	210848	2.367	ng	97
88) Benzo(b)fluoranthene	22.54	252	475825m	4.874	ng	
90) Benzo(a)pyrene	23.07	252	327119	3.685	ng	96
92) Benzo(g,h,i)perylene	25.86	276	223887	2.927	ng	97

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

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