

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP050520\
 Data File : BP002248.D
 Acq On : 05 May 2020 21:35
 Operator : CG/JU
 Sample : L2475-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FK-SF-03-009-A

Manual Integrations
 APPROVED

mohammad
 5/6/2020 6:44:42 PM

Quant Time: May 06 02:00:27 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP050420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 05 13:34:25 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	382638	20.00	ng	0.00
21) Naphthalene-d8	10.39	136	1675946	20.00	ng	0.00
39) Acenaphthene-d10	14.26	164	1024201	20.00	ng	0.00
64) Phenanthrene-d10	17.01	188	2193694	20.00	ng	0.00
76) Chrysene-d12	21.11	240	1622321	20.00	ng	0.00
87) Perylene-d12	23.32	264	1456436	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	2909190	125.43	ng	0.00
7) Phenol-d6	6.80	99	4048645	118.83	ng	0.01
23) Nitrobenzene-d5	8.76	82	2589060	79.78	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	1209838	111.60	ng	0.00
45) 2-Fluorobiphenyl	12.88	172	5053179	69.09	ng	0.00
79) Terphenyl-d14	19.60	244	5958447	90.01	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.72	163	410910	5.090	ng	99
71) Phenanthrene	17.05	178	689867	5.592	ng	99
75) Fluoranthene	19.03	202	1712425	12.099	ng	99
78) Pyrene	19.38	202	1604556	14.560	ng	99
81) Benzo(a)anthracene	21.09	228	744031	7.070	ng	96
83) Chrysene	21.15	228	741301m	7.173	ng	
86) Indeno(1,2,3-cd)pyrene	25.53	276	371309	2.640	ng	96
88) Benzo(b)fluoranthene	22.66	252	893262m	10.200	ng	
89) Benzo(k)fluoranthene	22.69	252	235203m	2.622	ng	
90) Benzo(a)pyrene	23.22	252	588192	6.976	ng	96
92) Benzo(g,h,i)perylene	26.20	276	388205	4.290	ng	98

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

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