

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP092019\
 Data File : BP000356.D
 Acq On : 20 Sep 2019 16:00
 Operator : HP/JU
 Sample : K4957-12
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-26-COMP

Manual Integrations
 APPROVED

mohammad
 9/24/2019 8:34:00 AM

Quant Time: Sep 20 23:49:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP091619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 19 14:44:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	172985	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	643431	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	364574	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	696128	20.00	ng	0.00
76) Chrysene-d12	13.99	240	625638	20.00	ng	0.00
87) Perylene-d12	15.47	264	728245	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	823717	82.21	ng	0.00
7) Phenol-d6	6.43	99	1056189	85.99	ng	0.00
23) Nitrobenzene-d5	7.35	82	577730	57.93	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	367610	101.66	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1249767	61.70	ng	0.00
79) Terphenyl-d14	12.93	244	1526895	51.20	ng	0.00
Target Compounds						
10) Phenol	6.44	94	49982	3.582	ng	99
50) Dimethylphthalate	9.54	163	199131	8.017	ng	99
71) Phenanthrene	11.35	178	297080	8.516	ng	98
72) Anthracene	11.40	178	88716	2.471	ng	99
75) Fluoranthene	12.55	202	644635	17.209	ng	95
78) Pyrene	12.78	202	613777	13.113	ng	99
81) Benzo(a)anthracene	13.98	228	358772	9.078	ng	97
83) Chrysene	14.02	228	309569	7.978	ng	97
86) Indeno(1,2,3-cd)pyrene	16.96	276	197048m	4.166	ng	
88) Benzo(b)fluoranthene	15.05	252	488650m	11.248	ng	
89) Benzo(k)fluoranthene	15.07	252	133561m	3.524	ng	
90) Benzo(a)pyrene	15.41	252	333615	8.439	ng	97
92) Benzo(a,h,i)perylene	17.40	276	184379	4.867	ng	95

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

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