

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090819\
 Data File : BP000126.D
 Acq On : 09 Sep 2019 06:25
 Operator : HP/JU
 Sample : K4736-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FK-GF-02-045-A

Manual Integrations
 APPROVED

mohammad
 9/10/2019 2:51:43 PM

Quant Time: Sep 09 08:51:22 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 30 17:47:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	424534	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	1576099	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	835815	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	1509674	20.00	ng	0.00
76) Chrysene-d12	14.10	240	1188497	20.00	ng	0.00
87) Perylene-d12	15.63	264	1059955	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	2392866	89.88	ng	0.00
7) Phenol-d6	6.52	99	3113852	92.65	ng	0.00
23) Nitrobenzene-d5	7.45	82	1691912	65.87	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	796348	113.04	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	3302581	65.99	ng	0.00
79) Terphenyl-d14	13.04	244	3567622	65.43	ng	0.00
Target Compounds						
10) Phenol	6.53	94	121652	3.258	ng	98
50) Dimethylphthalate	9.65	163	502293	8.819	ng	99
71) Phenanthrene	11.46	178	739248	9.825	ng	97
72) Anthracene	11.51	178	236608	3.183	ng	96
75) Fluoranthene	12.66	202	1164870	14.246	ng	96
78) Pyrene	12.89	202	1135469	12.987	ng	97
81) Benzo(a)anthracene	14.09	228	563134	7.312	ng	93
83) Chrysene	14.13	228	483460	6.637	ng	96
86) Indeno(1,2,3-cd)pyrene	17.18	276	215555m	2.361	ng	
88) Benzo(b)fluoranthene	15.17	252	584302m	9.025	ng	
89) Benzo(k)fluoranthene	15.20	252	144130m	2.433	ng	
90) Benzo(a)pyrene	15.56	252	405264	6.934	ng	97
92) Benzo(a,h,i)perylene	17.65	276	199463	3.448	ng	# 94

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

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