

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091819\
 Data File : BP000303.D
 Acq On : 18 Sep 2019 17:37
 Operator : HP/JU
 Sample : K4942-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-1

Manual Integrations
 APPROVED

Jagrut
 9/19/2019 11:07:09 AM

Quant Time: Sep 19 04:19:22 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP091619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 16 18:59:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	278964	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	1010535	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	549749	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	1019654	20.00	ng	0.00
76) Chrysene-d12	13.99	240	842667	20.00	ng	-0.01
87) Perylene-d12	15.47	264	920347	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	1569226	97.11	ng	0.00
7) Phenol-d6	6.43	99	1989769	100.45	ng	0.00
23) Nitrobenzene-d5	7.35	82	1125517	71.86	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	670240	122.92	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	2231066	73.04	ng	0.00
79) Terphenyl-d14	12.93	244	2537332	63.17	ng	0.00
Target Compounds						
10) Phenol	6.43	94	61986	2.754	ng	83
50) Dimethylphthalate	9.55	163	314168	8.388	ng	98
75) Fluoranthene	12.55	202	201609	3.674	ng	98
78) Pyrene	12.77	202	182375	2.893	ng	98
88) Benzo(b)fluoranthene	15.04	252	135595m	2.470	ng	

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

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