

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
 Data File : BP001389.D
 Acq On : 24 Dec 2019 16:35
 Operator : JU
 Sample : K6396-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-1-(5-7)

Manual Integrations
 APPROVED

mohammad
 12/26/2019 4:01:23 PM

Quant Time: Dec 25 02:33:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 16:04:49 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.28	152	55086	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	208876	20.00	ng	0.00
39) Acenaphthene-d10	13.18	164	126665	20.00	ng	0.00
64) Phenanthrene-d10	15.58	188	238744	20.00	ng	0.00
76) Chrysene-d12	19.22	240	158478	20.00	ng	0.00
87) Perylene-d12	20.99	264	150220	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.20	112	339324	99.56	ng	0.00
7) Phenol-d6	7.75	99	465689	104.71	ng	0.00
23) Nitrobenzene-d5	9.17	82	319693	58.77	ng	0.00
42) 2,4,6-Tribromophenol	14.47	330	143649	90.71	ng	0.00
45) 2-Fluorobiphenyl	12.10	172	530006	52.91	ng	0.00
79) Terphenyl-d14	17.86	244	478152	51.68	ng	0.00
Target Compounds						
10) Phenol	7.76	94	14895	2.931	ng	Qvalue # 79
50) Dimethylphthalate	12.76	163	26754	2.575	ng	99
71) Phenanthrene	15.61	178	40178	2.950	ng	99
75) Fluoranthene	17.33	202	45283	2.973	ng	98
78) Pyrene	17.63	202	37128	3.017	ng	# 94

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

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