

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011520\
 Data File : BP001620.D
 Acq On : 15 Jan 2020 21:21
 Operator : JU
 Sample : L1123-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-02-COMP

Manual Integrations
 APPROVED

mohammad
 1/16/2020 1:46:23 PM

Quant Time: Jan 16 07:29:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 14 21:52:04 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	32882	20.00	ng	0.00
21) Naphthalene-d8	10.40	136	116667	20.00	ng	-0.01
39) Acenaphthene-d10	14.28	164	79825	20.00	ng	0.00
64) Phenanthrene-d10	17.04	188	181026	20.00	ng	-0.01
76) Chrysene-d12	21.15	240	174970	20.00	ng	-0.01
87) Perylene-d12	23.37	264	187408	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	175280	104.42	ng	0.00
7) Phenol-d6	6.83	99	248684	92.70	ng	0.00
23) Nitrobenzene-d5	8.78	82	175676	65.80	ng	0.00
42) 2,4,6-Tribromophenol	15.78	330	97481	80.22	ng	-0.01
45) 2-Fluorobiphenyl	12.90	172	337577	62.18	ng	0.00
79) Terphenyl-d14	19.63	244	525691	54.95	ng	0.00
Target Compounds						
10) Phenol	6.85	94	7193	2.592	ng	86
50) Dimethylphthalate	13.75	163	37672	5.730	ng	99
75) Fluoranthene	19.07	202	31727	2.449	ng	94
78) Pyrene	19.42	202	31147	2.423	ng	98
83) Chrysene	21.18	228	25161	2.119	ng	96
88) Benzo(b)fluoranthene	22.71	252	35721m	3.025	ng	
90) Benzo(a)pyrene	23.28	252	25426	2.268	ng	99

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

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