

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100820\
 Data File : BP003582.D
 Acq On : 08 Oct 2020 21:13
 Operator : CG/JU
 Sample : L4224-01RE 5X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OR-02-100620RE

Manual Integrations
 APPROVED

mohammad
 10/12/2020 11:32:32 AM

Quant Time: Oct 09 02:00:19 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 08 15:55:58 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.481	152	60720	20.00	ng	0.00
21) Naphthalene-d8	10.258	136	234732	20.00	ng	0.00
39) Acenaphthene-d10	14.140	164	148265	20.00	ng	0.00
64) Phenanthrene-d10	16.898	188	322739	20.00	ng	0.00
76) Chrysene-d12	21.010	240	362861	20.00	ng	# 0.00
86) Perylene-d12	23.180	264	400038	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.111	112	29406	8.46	ng	0.01
7) Phenol-d6	6.717	99	33959	7.33	ng	0.03
23) Nitrobenzene-d5	8.652	82	26109	6.53	ng	0.02
42) 2,4,6-Tribromophenol	15.657	330	18334	8.11	ng	0.01
45) 2-Fluorobiphenyl	12.757	172	66350	6.54	ng	0.00
79) Terphenyl-d14	19.486	244	108555	6.23	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	16.945	178	60012	3.42	ng	98
75) Fluoranthene	18.934	202	98551	4.46	ng	98
78) Pyrene	19.286	202	90007	3.98	ng	99
83) Chrysene	21.045	228	48206	2.14	ng	# 92
88) Benzo(b)fluoranthene	22.533	252	56462m	2.24	ng	

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

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