

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010825.D
 Acq On : 25 Jun 2022 13:21
 Operator : CG/JU
 Sample : PB145664BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK664

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 06/27/2022
 Supervised By :mohammad ahmed 07/01/2022

Quant Time: Jun 27 09:30:43 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 24 22:42:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	402186	20.000	ng/u1	0.00
20) Naphthalene-d8	10.516	136	1621113	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.375	164	864336	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.139	188	1535499	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.251	240	1419877	20.000	ng/u1	# 0.00
88) Perylene-d12	23.621	264	1577719	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	77625	7.301	ng/uL	0.00
4) Pyridine-d5	3.616	84	1046087	36.963	ng/u1	0.01
7) Phenol-d5	6.904	99	1187920	37.312	ng/u1	0.01
9) Bis-(2-Chloroethyl)eth...	7.063	67	769586	36.961	ng/u1	0.00
11) 2-Chlorophenol-d4	7.257	132	1026050	38.845	ng/u1	0.00
15) 4-Methylphenol-d8	8.440	113	957515	38.045	ng/u1	0.00
21) Nitrobenzene-d5	8.881	128	466504	45.204	ng/u1	0.00
24) 2-Nitrophenol-d4	9.598	143	460711	53.960	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.139	165	796572	37.759	ng/u1	0.00
31) 4-Chloroaniline-d4	10.663	131	1303635	37.856	ng/u1	0.00
46) Dimethylphthalate-d6	13.798	166	2290521	38.686	ng/u1	0.00
49) Acenaphthylene-d8	14.069	160	2800552	37.061	ng/u1	0.00
54) 4-Nitrophenol-d4	14.598	143	323132m	34.941	ng/u1	0.02
60) Fluorene-d10	15.374	176	1894144	37.279	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.492	200	157671	30.993	ng/u1	0.00
73) Anthracene-d10	17.239	188	2709705	39.664	ng/u1	0.00
81) Pyrene-d10	19.486	212	2882125	37.258	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.468	264	3102231	39.791	ng/u1	0.00

Target Compounds Qvalue

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

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