

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030519\
 Data File : BF112858.D
 Acq On : 5 Mar 2019 17:56
 Operator : JU/SJ
 Sample : K1729-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RICH-REC-ROAD-STONE-10

Manual Integrations
 APPROVED

Sohil
 3/6/2019 12:24:25 PM

Quant Time: Mar 06 03:34:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 01 15:40:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	187726	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	733737	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	358494	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	724082	20.00	ng	0.00
76) Chrysene-d12	13.87	240	535876	20.00	ng	0.00
87) Perylene-d12	15.29	264	512760	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	164822	13.66	ng	0.00
7) Phenol-d6	6.37	99	827899	54.61	ng	0.00
23) Nitrobenzene-d5	7.30	82	763510	62.27	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	2507	0.66	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	1487833	64.53	ng	0.00
79) Terphenyl-d14	12.83	244	1593812	57.66	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.49	163	85486	3.234	ng	100
71) Phenanthrene	11.27	178	130148	3.613	ng	99
75) Fluoranthene	12.45	202	214223	5.550	ng	96
78) Pyrene	12.68	202	161435	3.573	ng	98
88) Benzo(b)fluoranthene	14.88	252	83594m	2.520	ng	

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

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